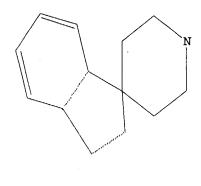
=> Uploading C:\Stnexp4 corrupted\QUERIES\10669606.str

STRUCTURE UPLOADED L12

=> d 112 L12 HAS NO ANSWERS

STR L12



G1 C, O, S, N

Structure attributes must be viewed using STN Express query preparation.

=> s 112 SAMPLE SEARCH INITIATED 12:06:26 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 481 TO ITERATE

100.0% PROCESSED 481 ITERATIONS 3 ANSWERS SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\* BATCH \*\*COMPLETE\*\*

8305 TO 10935 PROJECTED ITERATIONS:

PROJECTED ANSWERS: 3 TO 163

3 SEA SSS SAM L12 L13

=> s l12 ful

FULL SEARCH INITIATED 12:06:31 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 9456 TO ITERATE

20 ANSWERS 100.0% PROCESSED 9456 ITERATIONS SEARCH TIME: 00.00.01

L1420 SEA SSS FUL L12

=> file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 155.42 631.49

FILE 'CAPLUS' ENTERED AT 12:06:35 ON 09 JUL 2004

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FILE COVERS 1907 - 9 Jul 2004 VOL 141 ISS 3 FILE LAST UPDATED: 8 Jul 2004 (20040708/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 114

L15 1 L14

=> d abs bib hitstr

L15 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN GI

$$(R?)_p$$
 $L-B$ 
 $A$ 
 $A$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 
 $R^2$ 

III

```
The title compds. I and II [A = (substituted) 5- or 6-membered heteroarom.
AB
     ring or 5-, or 6-membered (hetero)cycloalkane ring; B =
      (hetero)cycloalkyl, (hetero)cycloalkenyl, or (hetero)aryl; L = alkylene;
     X, Y = a divalent linkage from the group consisting of alkylene,
     alkylene-OR3, alkylene-N(R3)COR4, alkylene-C(O)NR3R4, alkylene-N(R3)CO2R4,
     alkylene-N(R3)C(O)NR3R4, alkylene-C(O), O, C(O), N(R3), C(O)N(R3), S(O)k,
     and SO2N(R3); R1, R2 = H, alkyl, alkenyl, alkynyl, heteroalkyl, aryl,
     arylalkyl, NR6C(O)R5, C(O)R5, and NR5C(O)NR6; Rb = alkyl, aryl, OR7, C(O)R7, and C(O)NR7R8; R3, R4 = H, (hetero)alkyl, aryl, arylalkyl, etc.;
     R5, R6, R7, and R8 = H, alkyl, aryl, arylalkyl, etc.; p = 0-4; k = 0-2;
     with the proviso that X and Y are not both O, N(R3), S(O)k, or SO2N(R3)]
     were prepared as MCHR modulators in the treatment and/or prevention of
     endocrine, metabolic, cardiovascular, neurol., psychiatric,
     gastrointestinal, genitourinary and other disorders. For example,
     reaction of (2S)-2,3-dihydro-2-hydroxy-spiro[1H-indene-1,4'-piperidine]-1'-
     carboxylic acid 1,1-dimethylethyl ester (preparation given) with
     9-ethyl-9H-carbazole-3-carboxaldehyde yielded compound III. The prepared
     compds. were evaluated for MCHR2 modulation using a functional cell-based
     assay and most of them displayed potencies of 10 µM or less.
AN
     2004:290467 CAPLUS
DN
     140:303531
     Preparation of piperidinylmethyl carbazole derivatives as MCHR modulators
TI
     Chen, Xiaoqi; Dai, Kang; Fan, Pingchen; Fu, Ying; Li, Leping; Mihalic,
IN
     Jeffrey T.
     Tularik Inc., USA
PA
     PCT Int. Appl., 90 pp.
SO
     CODEN: PIXXD2
DT
     Patent
LA
     English
FAN.CNT 1
                                                 APPLICATION NO.
     PATENT NO.
                         KIND DATE
                                                  _____
      _____
     WO 2004028459
                        A2 20040408
                                                WO 2003-US30379 20030924
PΙ
          W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VC, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, PU, TJ
               KZ, MD, RU, TJ
          RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG,
               CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC,
               NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ,
               GW, ML, MR, NE, SN, TD, TG
PRAI US 2002-413335P
                                20020924
                          P
os
     MARPAT 140:303531
      676606-93-2P 676607-01-5P 676607-06-0P
TT
     RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic
     preparation); THU (Therapeutic use); BIOL (Biological study); PREP
      (Preparation); RACT (Reactant or reagent); USES (Uses)
          (preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)
RN
      676606-93-2 CAPLUS
      Acetamide, N-[(3R)-1'-(9H-carbazol-3-ylmethyl)-2,3-dihydrospiro[1H-indene-
CN
      1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)
```

Absolute stereochemistry.

RN 676607-01-5 CAPLUS CN Spiro[1H-indene-1,4'-piperidine], 1'-(9H-carbazol-3-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-06-0 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-3-acetic acid, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-, monosodium salt (9CI) (CA INDEX NAME)

Na

```
IT
     676606-89-6P 676606-90-9P 676606-92-1P
     676606-94-3P 676606-95-4P 676606-96-5P
     676606-97-6P 676606-98-7P 676607-00-4P
     676607-02-6P 676607-03-7P 676607-04-8P
     676607-05-9P 676607-07-1P 676607-08-2P
     676607-10-6P
     RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
        (preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)
RN
     676606-89-6 CAPLUS
     Spiro[1H-indene-1,4'-piperidin]-2-ol, 1'-[(9-ethyl-9H-carbazol-3-
CN
     yl)methyl]-2,3-dihydro-, (2S)- (9CI) (CA INDEX NAME)
```

Absolute stereochemistry.

RN 676606-90-9 CAPLUS
CN Acetamide, N-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydro-2-hydroxyspiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-92-1 CAPLUS
CN Spiro[1H-indene-1,4'-piperidin]-3-ol, 3-ethyl-1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydro- (9CI) (CA INDEX NAME)

RN 676606-94-3 CAPLUS
CN Urea, N-ethyl-N'-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-95-4 CAPLUS CN Carbamic acid, [1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydrospiro[1H-

indene-1,4'-piperidin]-3-yl]-, ethyl ester (9CI) (CA INDEX NAME)

RN 676606-96-5 CAPLUS
CN 4-Pyridineacetamide, N-[1'-[(9-ethyl-9H-carbazol-3-yl)methyl]-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-3-yl]- (9CI) (CA INDEX NAME)

RN 676606-97-6 CAPLUS CN Spiro[1H-indene-1,4'-piperidin]-2-ol, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-, (2R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 676606-98-7 CAPLUS
CN 9H-Carbazole-9-carboxamide, 3-[(2,3-dihydro-2-hydroxyspiro[1H-indene-1,4'-piperidin]-1'-yl)methyl]-N,N-dimethyl- (9CI) (CA INDEX NAME)

RN 676607-00-4 CAPLUS

CN 9H-Carbazole-9-carboxamide, 3-[[(3R)-3-(acetylamino)-2,3-dihydrospiro[1H-indene-1,4'-piperidin]-1'-yl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 676607-02-6 CAPLUS

CN 9H-Carbazole-9-carboxamide, 3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-03-7 CAPLUS

CN 9H-Carbazole, 9-(1-oxopropyl)-3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-04-8 CAPLUS

CN 9H-Carbazole-9-carboxamide, N,N-dimethyl-3-(spiro[1H-indene-1,4'-piperidin]-1'-ylmethyl)- (9CI) (CA INDEX NAME)

RN 676607-05-9 CAPLUS CN Spiro[1H-indene-1,4'-piperidine], 1'-[[9-(2-methoxyethyl)-9H-carbazol-3-yl]methyl]- (9CI) (CA INDEX NAME)

RN 676607-07-1 CAPLUS
CN Spiro[1H-indene-1,4'-piperidine]-3-acetamide, 1'-(9H-carbazol-3-ylmethyl)2,3-dihydro-N-3-pyridinyl- (9CI) (CA INDEX NAME)

RN 676607-08-2 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-3-acetamide, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-N-4-pyridinyl- (9CI) (CA INDEX NAME)

RN 676607-10-6 CAPLUS

CN Spiro[1H-indene-1,4'-piperidine]-3-carboxamide, 1'-(9H-carbazol-3-ylmethyl)-2,3-dihydro-N-(1-methylethyl)- (9CI) (CA INDEX NAME)

676607-27-5P IT

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of piperidinylmethyl carbazole derivs. as MCHR modulators)

676607-27-5 CAPLUS RN

Spiro[1H-indene-1,4'-piperidin]-3(2H)-one, 1'-[(9-ethyl-9H-carbazol-3-CNyl)methyl]- (9CI) (CA INDEX NAME)

=>